

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:sssptal204jxv

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * *
SESSION RESUMED IN FILE 'CAPLUS' AT 11:11:24 ON 02 MAY 2004
FILE 'CAPLUS' ENTERED AT 11:11:24 ON 02 MAY 2004
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COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	19.09	189.72

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-2.08	-2.08

=> file beilstein

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	19.96	190.59

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-2.08	-2.08

FILE 'BEILSTEIN' ENTERED AT 11:12:26 ON 02 MAY 2004
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FILE RELOADED ON OCTOBER 20, 2002
FILE LAST UPDATED ON MARCH 30,2004

FILE COVERS 1771 TO 2003.

*** FILE CONTAINS 8,932,479 SUBSTANCES ***

>>> PLEASE NOTE: Reaction data and substance data are stored in separate documents and can not be searched together in one query.

Reaction data for BEILSTEIN compounds may be displayed immediately with the display codes PRE (preparations) and REA (reactions). A substance answer set retrieved after the search for a chemical name, a molecular formula or a structure search for example can be restricted to compounds with available reaction information by concatenation with PRE/FA, REA/FA or more general with RX/FA. The BEILSTEIN Registry Number (BRN) is the link between a BEILSTEIN compound and belonging reactions. For more detailed reaction searches BRNs can be selected from substance answer sets and searched in the next step as reaction partner BRNs - Reactant (RX.RBRN) or Product BRN (RX.PBRN). After a search for reaction details substance documents associated with reactants or products may be retrieved by searching RX.PBRNs or RX.RBRNs as BRNs. <<<

>>> FOR SEARCHING PREPARATIONS SEE HELP PRE <<<

* PLEASE NOTE THAT THERE ARE NO FORMATS FREE OF COST. *
* SET NOTICE FEATURE: THE COST ESTIMATES CALCULATED FOR SET NOTICE *
* ARE BASED ON THE HIGHEST PRICE CATEGORY. THEREFORE; THESE *
* ESTIMATES MAY NOT REFLECT THE ACTUAL COSTS. *
* FOR PRICE INFORMATION SEE HELP COST *

=> d his

(FILE 'HOME' ENTERED AT 10:21:40 ON 02 MAY 2004)

FILE 'REGISTRY' ENTERED AT 10:21:48 ON 02 MAY 2004

L1 STRUCTURE UPLOADED

L2 1 S L1

L3 8 S L1 FULL

FILE 'REGISTRY' ENTERED AT 10:22:48 ON 02 MAY 2004

FILE 'CAPLUS' ENTERED AT 10:22:51 ON 02 MAY 2004

L4 3 S L3

FILE 'BEILSTEIN' ENTERED AT 11:12:26 ON 02 MAY 2004

=> s l1 full

FULL SEARCH INITIATED 11:12:33 FILE 'BEILSTEIN'

FULL SCREEN SEARCH COMPLETED - 498 TO ITERATE

100.0% PROCESSED 498 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.08

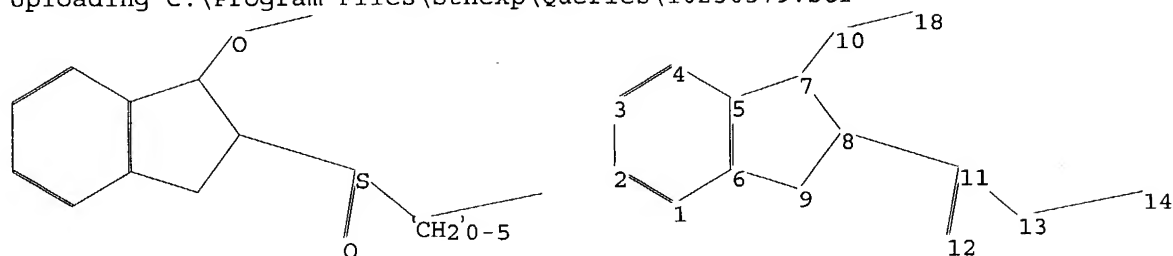
L5 1 SEA SSS FUL L1

=> d ide

L5 ANSWER 1 OF 1 BEILSTEIN COPYRIGHT 2004 BEILSTEIN MDL on STN

Beilstein Records (BRN):	2159196
Beilstein Pref. RN (BPR):	65495-98-9
CAS Reg. No. (RN):	65495-98-9
Chemical Name (CN):	1-Acetoxy-2-p-chlorophenylsulfinylindan
Autonom Name (AUN):	acetic acid 2-(4-chloro-benzenesulfinyl)-indan-1-yl ester
Molec. Formula (MF):	C17 H15 Cl O3 S
Molecular Weight (MW):	334.82
Lawson Number (LN):	6028, 5223, 1155
Compound Type (CTYPE):	isocyclic
Constitution ID (CONSID):	1965364
Tautomer ID (TAUTID):	2106089
Beilstein Citation (BSO):	5-06
Entry Date (DED):	1989/06/29
Update Date (DUPD):	1989/07/26

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 Uploading C:\Program Files\Stnexp\Queries\10230379.str

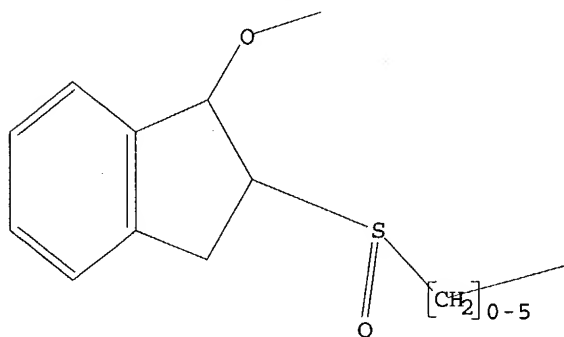


chain nodes :
 10 11 12 13
 ring nodes :
 1 2 3 4 5 6 7 8 9
 ring/chain nodes :
 14 18
 chain bonds :
 7-10 8-11 10-18 11-13 11-12 13-14
 ring bonds :
 1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9
 exact/norm bonds :
 7-10 8-11 10-18 11-12
 exact bonds :
 5-7 6-9 7-8 8-9 11-13 13-14
 normalized bonds :
 1-2 1-6 2-3 3-4 4-5 5-6
 isolated ring systems :
 containing 1 :

Match level :
 1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
 11:CLASS 12:CLASS 13:CLASS 14:CLASS 18:CLASS

L1 STRUCTURE UPLOADED

=> d
 L1 HAS NO ANSWERS
 L1 STR



Structure attributes must be viewed using STN Express query preparation.

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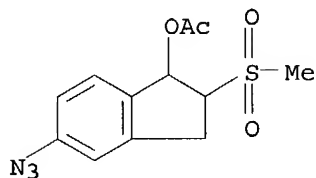
100.0% PROCESSED 96 ITERATIONS 1 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 1333 TO 2507
PROJECTED ANSWERS: 1 TO 80

L2 1 SEA SSS SAM L1

=> d scan

L2 1 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Inden-1-ol, 5-azido-2,3-dihydro-2-(methylsulfonyl)-, acetate (ester)
(9CI)
MF C12 H13 N3 O4 S



ALL ANSWERS HAVE BEEN SCANNED

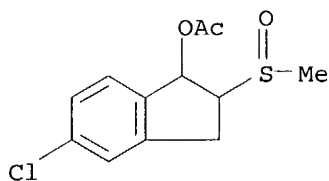
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FULL SCREEN SEARCH COMPLETED - 1860 TO ITERATE

100.0% PROCESSED 1860 ITERATIONS 8 ANSWERS
SEARCH TIME: 00.00.01

L3 8 SEA SSS FUL L1

=> d 1-8

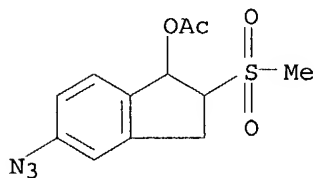
L3 ANSWER 1 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
RN 500770-94-5 REGISTRY
CN 1H-Inden-1-ol, 5-chloro-2,3-dihydro-2-(methylsulfinyl)-, acetate (9CI)
(CA INDEX NAME)
FS 3D CONCORD
MF C12 H13 Cl O3 S
SR CA
LC STN Files: CA, CAPLUS, USPAT2, USPATFULL



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

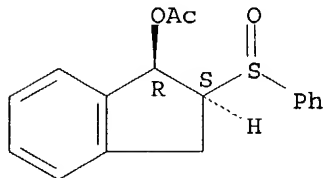
L3 ANSWER 2 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
RN 500770-93-4 REGISTRY
CN 1H-Inden-1-ol, 5-azido-2,3-dihydro-2-(methylsulfonyl)-, acetate (ester)
(9CI) (CA INDEX NAME)
FS 3D CONCORD
MF C12 H13 N3 O4 S
SR CA
LC STN Files: CA, CAPLUS, USPAT2, USPATFULL



1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 3 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
RN 500770-91-2 REGISTRY
CN 1H-Inden-1-ol, 2,3-dihydro-2-(phenylsulfinyl)-, acetate, (1R,2S)-rel-
(9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF C17 H16 O3 S
SR CA
LC STN Files: CA, CAPLUS, USPAT2, USPATFULL

Relative stereochemistry:



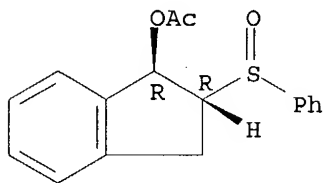
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1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 4 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
RN 500770-90-1 REGISTRY
CN 1H-Inden-1-ol, 2,3-dihydro-2-(phenylsulfinyl)-, acetate, (1R,2R)-rel-
(9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF C17 H16 O3 S
SR CA
LC STN Files: CA, CAPLUS, USPAT2, USPATFULL

Relative stereochemistry.

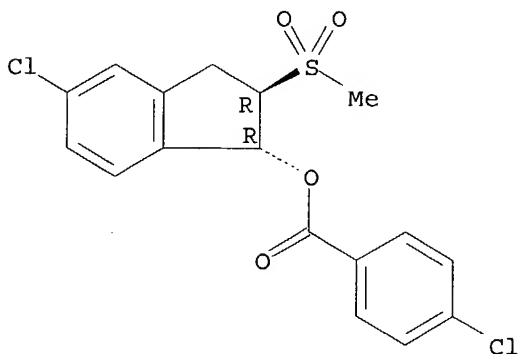


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 5 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
RN 500770-89-8 REGISTRY
CN Benzoic acid, 4-chloro-, (1R,2R)-5-chloro-2,3-dihydro-2-(methylsulfonyl)-
1H-inden-1-yl ester, rel- (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF C17 H14 Cl2 O4 S
SR CA
LC STN Files: CA, CAPLUS, USPAT2, USPATFULL

Relative stereochemistry.



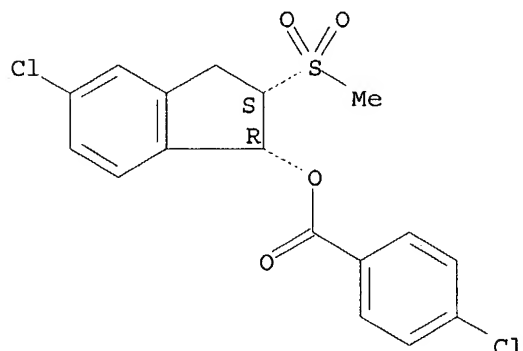
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2 REFERENCES IN FILE CA (1907 TO DATE)
2 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 6 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
RN 500770-88-7 REGISTRY

CN Benzoic acid, 4-chloro-, (1R,2S)-5-chloro-2,3-dihydro-2-(methylsulfonyl)-
 1H-inden-1-yl ester, rel- (9CI) (CA INDEX NAME)
 FS STEREOSEARCH
 MF C17 H14 Cl2 O4 S
 SR CA
 LC STN Files: CA, CAPLUS, USPAT2, USPATFULL

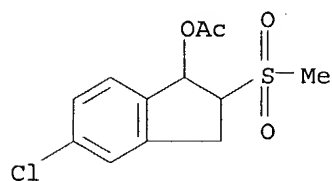
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2 REFERENCES IN FILE CA (1907 TO DATE)
 2 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 7 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 500770-87-6 REGISTRY
 CN 1H-Inden-1-ol, 5-chloro-2,3-dihydro-2-(methylsulfonyl)-, acetate (9CI)
 (CA INDEX NAME)
 FS 3D CONCORD
 MF C12 H13 Cl O4 S
 SR CA
 LC STN Files: CA, CAPLUS, USPAT2, USPATFULL

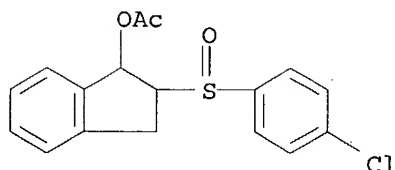


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2 REFERENCES IN FILE CA (1907 TO DATE)
 2 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 8 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 65495-98-9 REGISTRY
 CN 1H-Inden-1-ol, 2-[(4-chlorophenyl)sulfinyl]-2,3-dihydro-, acetate (9CI)
 (CA INDEX NAME)
 FS 3D CONCORD
 MF C17 H15 Cl O3 S
 LC STN Files: BEILSTEIN*, CA, CAPLUS, CASREACT

(*File contains numerically searchable property data)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> file reg

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
170.00	170.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 10:22:48 ON 02 MAY 2004
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Property values tagged with IC are from the ZIC/VINITI data file
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STRUCTURE FILE UPDATES: 30 APR 2004 HIGHEST RN 678535-01-8
DICTIONARY FILE UPDATES: 30 APR 2004 HIGHEST RN 678535-01-8

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
0.42	170.63

FULL ESTIMATED COST

FILE 'CAPLUS' ENTERED AT 10:22:51 ON 02 MAY 2004
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FILE COVERS 1907 - 2 May 2004 VOL 140 ISS 19
FILE LAST UPDATED: 30 Apr 2004 (20040430/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l3

L4 3 L3

=> d ibib abs hitstr 1-3

L4 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN

ACCESSION NUMBER: 2003:202616 CAPLUS

DOCUMENT NUMBER: 138:221363

TITLE: Preparation of indan-1-ol acetates and related compounds as appetite depressants

INVENTOR(S): Jaehne, Gerhard; Krone, Volker; Bickel, Martin; Gossel, Matthias

PATENT ASSIGNEE(S): Aventis Pharma Deutschland G.m.b.H., Germany

SOURCE: PCT Int. Appl., 54 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: German

FAMILY ACC. NUM. COUNT: 1

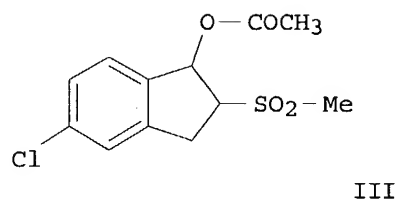
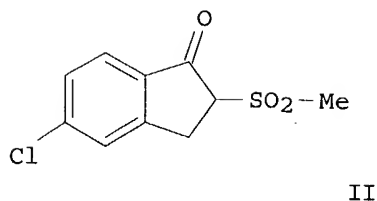
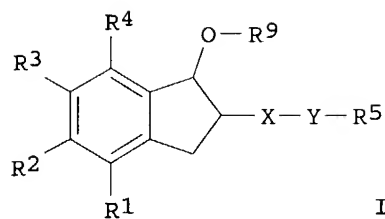
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003020694	A1	20030313	WO 2002-EP9202	20020817
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM			
RW:	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
DE 10142662	A1	20030327	DE 2001-10142662	20010831
US 2003105331	A1	20030605	US 2002-231365	20020830
US 6717008	B2	20040406		

PRIORITY APPLN. INFO.: DE 2001-10142662 A 20010831

OTHER SOURCE(S): MARPAT 138:221363

GI



AB Title compds. I [R1, R2, R3, R4 = H, halo, CN, etc.; X = S, SO, SO2; Y = (CH2)p; p = 0-3; R5 = CF3, alkyl, cycloalkyl; R9 = substituted alkyl or cycloalkyl, e.g. F, CO, CO2, etc.] and their pharmaceutically acceptable salts were prepared For example, NaBH4 mediated reduction of indanone II, e.g.,

prepared from 2-bromo-5-chloroindan-1-one in 2-steps, followed by O-acetylation provided indanol acetate III. In milk consumption studies with female NMRI mice, indanol III exhibited very good anorectic effects, i.e., 81% decrease in milk consumption verses control.

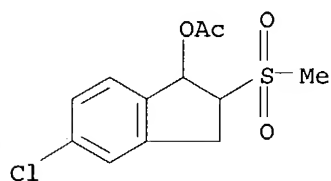
IT 500770-87-6P 500770-88-7P 500770-89-8P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(drug candidate; preparation of indanol acetates and related compds. as appetite depressants)

RN 500770-87-6 CAPLUS

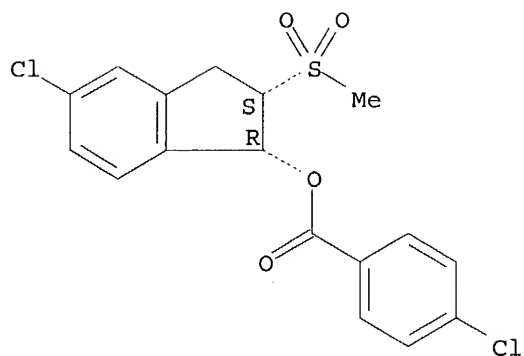
CN 1H-Inden-1-ol, 5-chloro-2,3-dihydro-2-(methylsulfonyl)-, acetate (9CI)
(CA INDEX NAME)



RN 500770-88-7 CAPLUS

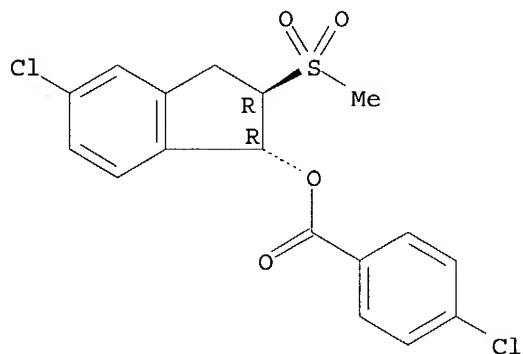
CN Benzoic acid, 4-chloro-, (1R,2S)-5-chloro-2,3-dihydro-2-(methylsulfonyl)-1H-inden-1-yl ester, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RN 500770-89-8 CAPLUS
 CN Benzoic acid, 4-chloro-, (1R,2R)-5-chloro-2,3-dihydro-2-(methanesulfonyl)-1H-inden-1-yl ester, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN

ACCESSION NUMBER: 2003:202409 CAPLUS

DOCUMENT NUMBER: 138:226750

TITLE: Use of C2-substituted indan-1-ol derivatives in antiobesity drugs

INVENTOR(S): Jaehne, Gerhard; Krone, Volker; Bickel, Martin; Gossel, Matthias

PATENT ASSIGNEE(S): Aventis Pharma Deutschland G.m.b.H., Germany

SOURCE: PCT Int. Appl., 54 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: German

FAMILY ACC. NUM. COUNT: 1

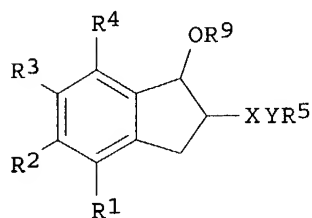
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003020199	A1	20030313	WO 2002-EP9199	20020817
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UA, UG, UZ, VN, YU, ZA, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
 RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, BG,
 CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL,
 PT, SE, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR,
 NE, SN, TD, TG

DE 10142660 A1 20030320 DE 2001-10142660 20010831
 US 2003134881 A1 20030717 US 2002-230379 20020829
 US 6667345 B2 20031223

PRIORITY APPLN. INFO.: DE 2001-10142660 A 20010831
 OTHER SOURCE(S): MARPAT 138:226750
 GI



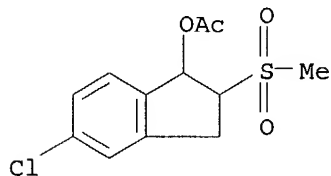
AB The invention relates to the use of C2-substituted indan-1-ol systems, and to the physiol. tolerable salts and the physiol. functional derivs. of the same, for producing medicaments used to reduce the weight of mammals, and for the prophylaxis or the treatment of obesity. The invention also relates to the use of compds. of formula (I), wherein the radicals have the cited designations, and to the physiol. tolerable salts and the physiol. functional derivs. of the same, for producing a medicament for the prophylaxis or the treatment of obesity. The antiobesity drugs can be combined with other active ingredients, e.g. cathine, phenylpropanolamine, amfepramone, mefenorex. Capsules, tablets, emulsions, dragees and suppositories are prepared containing the indan-1-ol derivative antiobesity drugs.

IT 500770-87-6P 500770-88-7P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
 (use of C2-substituted indan-1-ol derivs. in antiobesity drugs)

RN 500770-87-6 CAPLUS

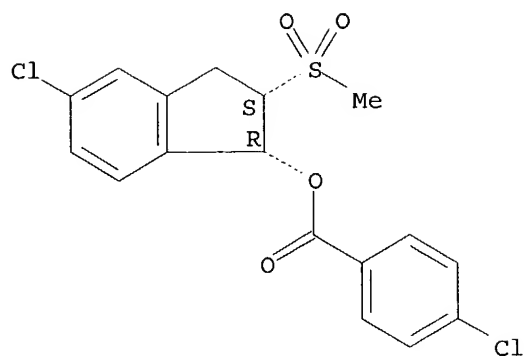
CN 1H-Inden-1-ol, 5-chloro-2,3-dihydro-2-(methylsulfonyl)-, acetate (9CI)
 (CA INDEX NAME)



RN 500770-88-7 CAPLUS

CN Benzoic acid, 4-chloro-, (1R,2S)-5-chloro-2,3-dihydro-2-(methylsulfonyl)-1H-inden-1-yl ester, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



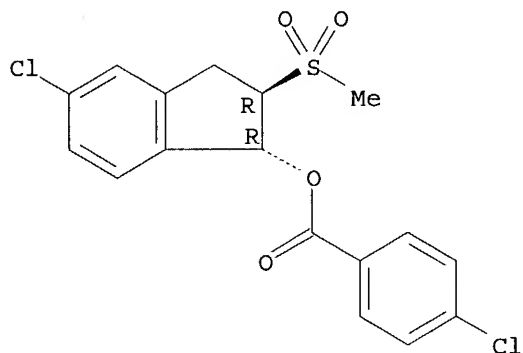
IT 500770-89-8 500770-90-1 500770-91-2
500770-93-4 500770-94-5

RL: THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(use of C2-substituted indan-1-ol derivs. in antiobesity drugs)

RN 500770-89-8 CAPLUS

CN Benzoic acid, 4-chloro-, (1R,2R)-5-chloro-2,3-dihydro-2-(methylsulfonyl)-
1H-inden-1-yl ester, rel- (9CI) (CA INDEX NAME)

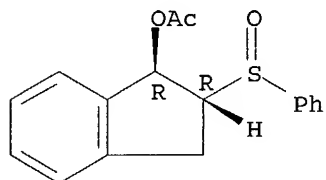
Relative stereochemistry.



RN 500770-90-1 CAPLUS

CN 1H-Inden-1-ol, 2,3-dihydro-2-(phenylsulfinyl)-, acetate, (1R,2R)-rel-
(9CI) (CA INDEX NAME)

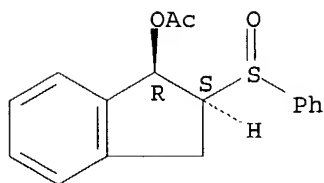
Relative stereochemistry.



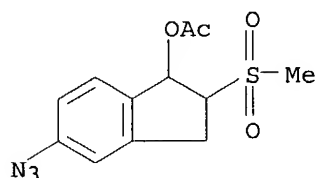
RN 500770-91-2 CAPLUS

CN 1H-Inden-1-ol, 2,3-dihydro-2-(phenylsulfinyl)-, acetate, (1R,2S)-rel-
(9CI) (CA INDEX NAME)

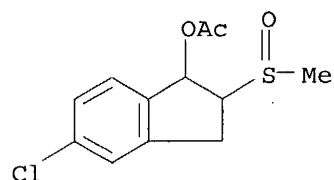
Relative stereochemistry.



RN 500770-93-4 CAPLUS
 CN 1H-Inden-1-ol, 5-azido-2,3-dihydro-2-(methylsulfonyl)-, acetate (ester)
 (9CI) (CA INDEX NAME)



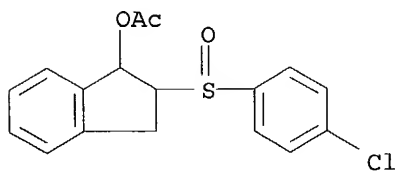
RN 500770-94-5 CAPLUS
 CN 1H-Inden-1-ol, 5-chloro-2,3-dihydro-2-(methylsulfinyl)-, acetate (9CI)
 (CA INDEX NAME)

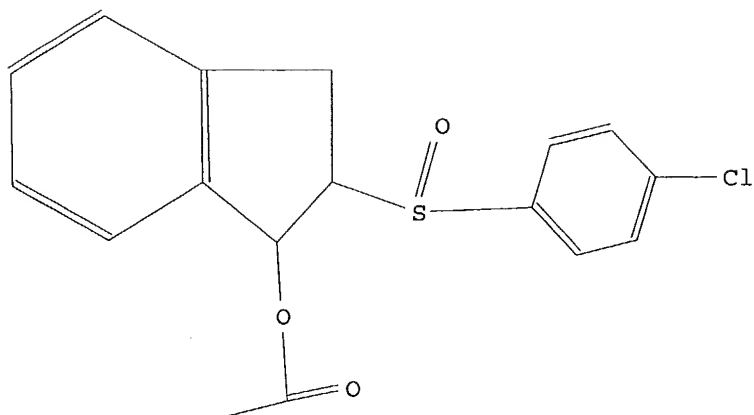


REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS
 RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN
 ACCESSION NUMBER: 1978:190435 CAPLUS
 DOCUMENT NUMBER: 88:190435
 TITLE: Thiol-olefin cooxidation reaction. 6. A new
 convenient route to 1-substituted indenenes. Indenone
 as dienophile in Diels-Alder reactions
 AUTHOR(S): Szmant, H. Harry; Nanjundiah, Raghunath
 CORPORATE SOURCE: Dep. Chem. Chem. Eng., Univ. Detroit, Detroit, MI, USA
 SOURCE: Journal of Organic Chemistry (1978), 43(9), 1835-7
 CODEN: JOCEAH; ISSN: 0022-3263
 DOCUMENT TYPE: Journal
 LANGUAGE: English
 OTHER SOURCE(S): CASREACT 88:190435
 AB 2-(4-Chlorophenylsulfinyl)-1-indanone was decomposed in refluxing toluene to
 give indenone which was trapped by cyclopentadiene,
 hexachlorocyclopentadiene, and anthracene to give the resp. Diels-Alder
 adducts.
 IT 65495-98-9P
 RL: SPN (Synthetic preparation); PREP (Preparation)
 (preparation of)
 RN 65495-98-9 CAPLUS

CN 1H-Inden-1-ol, 2-[(4-chlorophenyl)sulfinyl]-2,3-dihydro-, acetate (9CI)
(CA INDEX NAME)





Field Availability:

Code	Name	Occurrence
BRN	Beilstein Records	1
BPR	Beilstein Preferred RN	1
RN	CAS Registry Number	1
CN	Chemical Name	1
AUN	Autonomname	1
MF	Molecular Formula	1
FW	Formular Weight	1
LN	Lawson Number	3
CTYPE	Compound Type	1
CONSID	Constitution ID	1
TAUTID	Tautomer ID	1
BSO	Beilstein Citation	1
ED	Entry Date	1
UPD	Update Date	1
MP	Melting Point	1
NMR	Nuclear Magnetic Resonance	1

This substance also occurs in Reaction Documents:

Code	Name	Occurrence
RX	Reaction Documents	1
RXPRO	Substance is Reaction Product	1

=> d rxpro

L5 ANSWER 1 OF 1 BEILSTEIN COPYRIGHT 2004 BEILSTEIN MDL on STN

Reaction:

RX

Reaction ID (.ID): 7329680
 Product BRN (.PBRN): 2159196
 Product (.PRO): 1-Acetoxy-2-p-chlorophenylsulfinylindan
 No. of React. Details (.NVAR): 1

Reaction Details:

RX

Reaction RID (.RID): 7329680.1
 Reaction Classification (.CL): Preparation (half reaction)
 Reference(s):

1. Szmant; Nanjundiah, J.Org.Chem., CODEN: JOCEAH, 43, <1978>, 1835